

EAST Search History

Ref #	Hits	Search Query	DBs	Default Operator	Plurals	Time Stamp
L1	519538	phenyl	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/01 08:56
L2	2354	hexenoic	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/01 08:56
L3	39	L2 near5 L1	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/01 09:56
L4	176	(562/491).CCLS.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	OFF	2006/12/01 09:02
L5	241	(562/495).CCLS.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	OFF	2006/12/01 09:02
L6	590	(514/559).CCLS.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	OFF	2006/12/01 09:03
L7	1354	(514/562).CCLS.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	OFF	2006/12/01 09:03
L8	571	(514/564).CCLS.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	OFF	2006/12/01 09:03
L9	1066	(514/570).CCLS.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	OFF	2006/12/01 09:03
L10	319	(514/571).CCLS.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	OFF	2006/12/01 09:03
L11	3931	I4 or I5 or I6 or I7 or I8 or I9 or I10	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/01 09:04

EAST Search History

L12	1	I3 and I11	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/01 09:04
L13	150554	cyclohexyl	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/01 09:56
L14	16	I3 and I13	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/01 09:56
L15	11	("4513005").URPN.	USPAT	OR	ON	2006/12/01 09:56

CA SUBSCRIBER PRICE

ENTRY	SESSION
0.00	-2.25

SESSION WILL BE HELD FOR 120 MINUTES
STN INTERNATIONAL SESSION SUSPENDED AT 06:41:49 ON 01 DEC 2006

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:SSSPTA1623PAZ

PASSWORD:

***** RECONNECTED TO STN INTERNATIONAL *****
SESSION RESUMED IN FILE 'REGISTRY' AT 06:47:45 ON 01 DEC 2006
FILE 'REGISTRY' ENTERED AT 06:47:45 ON 01 DEC 2006
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COST IN U.S. DOLLARS	SINCE FILE	TOTAL
FULL ESTIMATED COST	ENTRY	SESSION
	7.84	243.93
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	0.00	-2.25

=> e 6-phenyl-2-hexynoic acid/cn

E1	1	6-PHENYL-2-HEXYN-1-OL/CN
E2	1	6-PHENYL-2-HEXYNE/CN
E3	1 -->	6-PHENYL-2-HEXYNOIC ACID/CN
E4	1	6-PHENYL-2-HYDROXY-N-(2-(3-METHOXY-4-((2-PROPYNYL) OXY) PHENYL) ETHYL) HEXANAMIDE/CN
E5	1	6-PHENYL-2-INDOLINONE/CN
E6	1	6-PHENYL-2-MERCAPTOTHIAZOLO(4,5-B) PYRIDINE/CN
E7	1	6-PHENYL-2-METHYL-5,6-DIHYDROOXAZINE/CN
E8	1	6-PHENYL-2-METHYLTHIO-THIAZOLO(4,5-B) PYRIDINE/CN
E9	1	6-PHENYL-2-NAPHTHOL/CN
E10	1	6-PHENYL-2-OCTANONE/CN
E11	1	6-PHENYL-2-OXINDOLE/CN
E12	1	6-PHENYL-2-PHENYLAMINO-4H-THIENO(3,2-D)(1,3) OXAZIN-4-ONE/CN

=> e3

L16 1 "6-PHENYL-2-HEXYNOIC ACID"/CN

=> d 116

L16 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2006 ACS on STN
RN 235087-20-4 REGISTRY
ED Entered STN: 26 Aug 1999
CN 2-Hexynoic acid, 6-phenyl- (9CI) (CA INDEX NAME)
OTHER NAMES:
CN 6-Phenyl-2-hexynoic acid
MF C12 H12 O2
SR CA
LC STN Files: CA, CAPLUS, CASREACT

HO₂C-C≡C-(CH₂)₃-Ph

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

2 REFERENCES IN FILE CA (1907 TO DATE)
2 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> file caplus

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
17.58	253.67

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION
0.00	-2.25

CA SUBSCRIBER PRICE

FILE 'CAPLUS' ENTERED AT 06:51:49 ON 01 DEC 2006

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FILE COVERS 1907 - 1 Dec 2006 VOL 145 ISS 23

FILE LAST UPDATED: 29 Nov 2006 (20061129/ED)

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<http://www.cas.org/infopolicy.html>

=> l16

L17 2 L16

=> d l17 1-2 ti fib abs

'FIB' IS NOT A VALID FORMAT FOR FILE 'CAPLUS'

The following are valid formats:

ABS ----- GI and AB
ALL ----- BIB, AB, IND, RE
APPS ----- AI, PRAI
BIB ----- AN, plus Bibliographic Data and PI table (default)
CAN ----- List of CA abstract numbers without answer numbers
CBIB ----- AN, plus Compressed Bibliographic Data
CLASS ----- IPC, NCL, ECLA, FTERM
DALL ----- ALL, delimited (end of each field identified)
DMAX ----- MAX, delimited for post-processing
FAM ----- AN, PI and PRAI in table, plus Patent Family data
FBIB ----- AN, BIB, plus Patent FAM
IND ----- Indexing data
IPC ----- International Patent Classifications
MAX ----- ALL, plus Patent FAM, RE
PATS ----- PI, SO

SAM ----- CC, SX, TI, ST, IT
 SCAN ----- CC, SX, TI, ST, IT (random display, no answer numbers;
 SCAN must be entered on the same line as the DISPLAY,
 e.g., D SCAN or DISPLAY SCAN)
 STD ----- BIB, CLASS

 IABS ----- ABS, indented with text labels
 IALL ----- ALL, indented with text labels
 IBIB ----- BIB, indented with text labels
 IMAX ----- MAX, indented with text labels
 ISTD ----- STD, indented with text labels

 OBIB ----- AN, plus Bibliographic Data (original)
 OIBIB ----- OBIB, indented with text labels

 SBIB ----- BIB, no citations
 SIBIB ----- IBIB, no citations

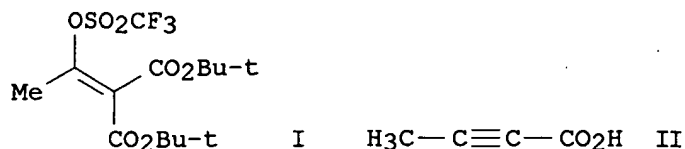
 HIT ----- Fields containing hit terms
 HITIND ----- IC, ICA, ICI, NCL, CC and index field (ST and IT)
 containing hit terms
 HITRN ----- HIT RN and its text modification
 HITSTR ----- HIT RN, its text modification, its CA index name, and
 its structure diagram
 HITSEQ ----- HIT RN, its text modification, its CA index name, its
 structure diagram, plus NTE and SEQ fields
 FHITSTR ----- First HIT RN, its text modification, its CA index name, and
 its structure diagram
 FHITSEQ ----- First HIT RN, its text modification, its CA index name, its
 structure diagram, plus NTE and SEQ fields
 KWIC ----- Hit term plus 20 words on either side
 OCC ----- Number of occurrence of hit term and field in which it occurs

To display a particular field or fields, enter the display field codes. For a list of the display field codes, enter HELP DFIELDS at an arrow prompt (=>). Examples of formats include: TI; TI,AU; BIB,ST; TI,IND; TI,SO. You may specify the format fields in any order and the information will be displayed in the same order as the format specification.

All of the formats (except for SAM, SCAN, HIT, HITIND, HITRN, HITSTR, FHITSTR, HITSEQ, FHITSEQ, KWIC, and OCC) may be used with DISPLAY ACC to view a specified Accession Number.
 ENTER DISPLAY FORMAT (BIB):end

=> d 117 1-2 ti fbib abs

L17 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2006 ACS on STN
 TI Decarboxylative elimination of enol triflates as a general synthesis of
 acetylenes
 AN 2004:380047 CAPLUS
 DN 141:140115
 TI Decarboxylative elimination of enol triflates as a general synthesis of
 acetylenes
 AU Fleming, Ian; Ramarao, Chandrashekar
 CS Department of Chemistry, Cambridge, CB2 1EW, UK
 SO Organic & Biomolecular Chemistry (2004), 2(10), 1504-1510
 CODEN: OBCRAK; ISSN: 1477-0520
 PB Royal Society of Chemistry
 DT Journal
 LA English
 OS CASREACT 141:140115
 GI



AB The enol trifluoromethanesulfonates of tert-Bu β -keto diesters and β -keto esters can be hydrolyzed to the corresponding carboxylic acids by dissoln. in trifluoroacetic acid. The dicarboxylic acids undergo mild decarboxylative elimination to give acetylenic acids in aqueous sodium bicarbonate solution at room temperature. Similarly, monocarboxylic acids give terminal and mid-chain acetylenes by refluxing in acetone with potassium carbonate. One of the substituents on the acetylenes can be Me, primary alkyl, secondary alkyl or ethynyl, and the other can be a carboxylic acid, hydrogen or primary alkyl, but the enol trifluoromethanesulfonates could not be prepared when one of the substituents was tert-Bu, nor when both substituents on the precursor to the acetylene were secondary alkyl. For example, reaction of trifluoromethanesulfonic acid anhydride with (acetyl)propanedioic acid bis(1,1-dimethylethyl) ester gave a desired enol triflate, [1-[[[(trifluoromethyl)sulfonyl]oxy]ethylidene]propanedioic acid bis(1,1-dimethylethyl) ester (I). Saponification of I gave the diacid, [1-[[[(trifluoromethyl)sulfonyl]oxy]ethylidene]propanedioic acid. Decarboxylation of the acid gave 2-butyric acid (II).

RE.CNT 46 THERE ARE 46 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L17 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2006 ACS on STN

TI Decarboxylative elimination of enol triflates as a general synthesis of acetylenes

AN 1999:351377 CAPLUS

DN 131:144304

TI Decarboxylative elimination of enol triflates as a general synthesis of acetylenes

AU Fleming, Ian; Ramarao, Chandrashekar

CS Department of Chemistry, Cambridge, CB2 1EW, UK

SO Chemical Communications (Cambridge) (1999), (12), 1113-1114

CODEN: CHCOFS; ISSN: 1359-7345

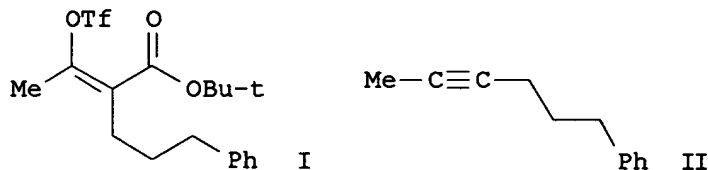
PB Royal Society of Chemistry

DT Journal

LA English

OS CASREACT 131:144304

GI



AB Decarboxylative elimination of a range of enol triflates, e.g. I, of β -keto esters gives acetylenes, e.g. II.

RE.CNT 17 THERE ARE 17 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> logoff hold

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
15.60	269.27

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION
-1.50	-3.75

CA SUBSCRIBER PRICE

SESSION WILL BE HELD FOR 120 MINUTES

STN INTERNATIONAL SESSION SUSPENDED AT 07:05:10 ON 01 DEC 2006

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:SSSPTA1623PAZ

PASSWORD:

* * * * * RECONNECTED TO STN INTERNATIONAL * * * * *
SESSION RESUMED IN FILE 'CAPLUS' AT 07:17:01 ON 01 DEC 2006
FILE 'CAPLUS' ENTERED AT 07:17:01 ON 01 DEC 2006
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COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
15.60	269.27

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION
-1.50	-3.75

CA SUBSCRIBER PRICE

=> file reg

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
15.60	269.27

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION
-1.50	-3.75

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Property values tagged with IC are from the ZIC/VINITI data file
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STRUCTURE FILE UPDATES: 29 NOV 2006 HIGHEST RN 914337-13-6

DICTIONARY FILE UPDATES: 29 NOV 2006 HIGHEST RN 914337-13-6

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH June 30, 2006

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and

predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

<http://www.cas.org/ONLINE/UG/regprops.html>

=> e 7-phenyl-2-heptynoic acid/cn

E1	1	7-PHENYL-2-ANILINO-1-PHENYL-1,8-NAPHTHYRIDIN-4(1H)-ONE/CN
E2	1	7-PHENYL-2-HEPTANONE/CN
E3	0 -->	7-PHENYL-2-HEPTYNOIC ACID/CN
E4	1	7-PHENYL-2-NAPHTHALENOL/CN
E5	1	7-PHENYL-2-NAPHTHOL/CN
E6	1	7-PHENYL-2-OCTANONE/CN
E7	1	7-PHENYL-2-OXA-7-AZABICYCLO(3.2.0)HEPTAN-6-ONE/CN
E8	1	7-PHENYL-2-OXEPANONE/CN
E9	1	7-PHENYL-3,4-DIHYDRO-1(2H)-NAPHTHALENONE/CN
E10	1	7-PHENYL-3,6-DIOXAHEPTYL P-TOLUENESULFONATE/CN
E11	1	7-PHENYL-3,6-DIOXAHEPTYL TOSYLATE/CN
E12	1	7-PHENYL-3-(2-(4-PYRIDYL)-1,3-THIAZOL-4-YL)-1,2,3,4-TETRAHYDROQUINAZOLIN-2-ONE/CN

=> e 6-phenylhexenoic acid/cn

E1	1	6-PHENYLHEXANOYLHYDROXAMIC ACID/CN
E2	1	6-PHENYLHEXENE/CN
E3	0 -->	6-PHENYLHEXENOIC ACID/CN
E4	1	6-PHENYLHEXYL 4-HYDROXYBENZOATE/CN
E5	1	6-PHENYLHEXYL ACETATE/CN
E6	1	6-PHENYLHEXYL ACRYLATE-METHYL METHACRYLATE-POLYETHYLENE GLYCOL METHYL ETHER ACRYLATE GRAFT COPOLYMER/CN
E7	1	6-PHENYLHEXYL BROMIDE/CN
E8	1	6-PHENYLHEXYL IODIDE/CN
E9	1	6-PHENYLHEXYL ISOTHIOCYANATE/CN
E10	1	6-PHENYLHEXYL MESYLATE/CN
E11	1	6-PHENYLHEXYLAMINE/CN
E12	1	6-PHENYLHEXYLMAGNESIUM BROMIDE/CN

=> e 6-phenyl-2-hexenoic acid/cn

E1	1	6-PHENYL-2-HEXEN-5-YNE/CN
E2	1	6-PHENYL-2-HEXENE/CN
E3	0 -->	6-PHENYL-2-HEXENOIC ACID/CN
E4	1	6-PHENYL-2-HEXYN-1-OL/CN
E5	1	6-PHENYL-2-HEXYNE/CN
E6	1	6-PHENYL-2-HEXYNOIC ACID/CN
E7	1	6-PHENYL-2-HYDROXY-N-(2-(3-METHOXY-4-((2-PROPYNYL)OXY)PHENYL)ETHYL)HEXANAMIDE/CN
E8	1	6-PHENYL-2-INDOLINONE/CN
E9	1	6-PHENYL-2-MERCAPTOTHIAZOLO(4,5-B)PYRIDINE/CN
E10	1	6-PHENYL-2-METHYL-5,6-DIHYDROOXAZINE/CN
E11	1	6-PHENYL-2-METHYLTHIO-THIAZOLO(4,5-B)PYRIDINE/CN
E12	1	6-PHENYL-2-NAPHTHOL/CN

=> e6

L18 1 "6-PHENYL-2-HEXYNOIC ACID"/CN

=> e 6-phenoxy-2-hexenoic acid/cn

E1	1	6-PHENOXY-2-((4-PHENOXYPHENYL)AMINO)-4H-THIENO(2,3-D)(1,3)OXAZIN-4-ONE/CN
E2	1	6-PHENOXY-2-BENZOFURANCARBOXYLIC ACID/CN
E3	0 -->	6-PHENOXY-2-HEXENOIC ACID/CN
E4	1	6-PHENOXY-2-HYDROXYHEXANOIC ACID/CN
E5	1	6-PHENOXY-2-OXOHEXANOIC ACID/CN
E6	1	6-PHENOXY-2-PYRIDINECARBOTHIOAMIDE/CN
E7	1	6-PHENOXY-2-PYRIDINECARBOXALDEHYDE/CN

E8	1	6-PHENOXY-2-PYRIDINECARBOXYLIC ACID/CN
E9	1	6-PHENOXY-2-PYRIDINEMETHANETHIOL/CN
E10	1	6-PHENOXY-2-PYRIDINEMETHANOL/CN
E11	1	6-PHENOXY-3-PYRIDINAMINE/CN
E12	1	6-PHENOXY-3-PYRIDINAMINE HYDROCHLORIDE/CN

=> d cost

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
CONNECT CHARGES	4.18	50.01
NETWORK CHARGES	0.66	7.74
SEARCH CHARGES	4.76	199.82
DISPLAY CHARGES	0.00	21.30
FULL ESTIMATED COST	9.60	278.87

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-3.75

IN FILE 'REGISTRY' AT 07:23:45 ON 01 DEC 2006

=> logoff hold

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	9.60	278.87
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-3.75

SESSION WILL BE HELD FOR 120 MINUTES
STN INTERNATIONAL SESSION SUSPENDED AT 07:23:52 ON 01 DEC 2006

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:SSSPTA1623PAZ

PASSWORD:

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SESSION RESUMED IN FILE 'REGISTRY' AT 09:09:42 ON 01 DEC 2006
FILE 'REGISTRY' ENTERED AT 09:09:42 ON 01 DEC 2006
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COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	9.60	278.87
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-3.75

=> logoff hold

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	9.60	278.87

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE

TOTAL

ENTRY

SESSION

CA SUBSCRIBER PRICE

0.00

-3.75

SESSION WILL BE HELD FOR 120 MINUTES

STN INTERNATIONAL SESSION SUSPENDED AT 09:09:50 ON 01 DEC 2006